

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: sssptal611txm

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * * * * * * * * Welcome to STN International * * * * * * * * * * * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 Apr 08 "Ask CAS" for self-help around the clock
NEWS 3 Jun 03 New e-mail delivery for search results now available
NEWS 4 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN
NEWS 5 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE)
now available on STN
NEWS 6 Aug 26 Sequence searching in REGISTRY enhanced
NEWS 7 Sep 03 JAPIO has been reloaded and enhanced
NEWS 8 Sep 16 Experimental properties added to the REGISTRY file
NEWS 9 Sep 16 CA Section Thesaurus available in CAPLUS and CA
NEWS 10 Oct 01 CASREACT Enriched with Reactions from 1907 to 1985
NEWS 11 Oct 24 BEILSTEIN adds new search fields
NEWS 12 Oct 24 Nutraceuticals International (NUTRACEUT) now available on STN
NEWS 13 Nov 18 DKILIT has been renamed APOLLIT
NEWS 14 Nov 25 More calculated properties added to REGISTRY
NEWS 15 Dec 04 CSA files on STN
NEWS 16 Dec 17 PCTFULL now covers WP/PCT Applications from 1978 to date
NEWS 17 Dec 17 TOXCENTER enhanced with additional content
NEWS 18 Dec 17 Adis Clinical Trials Insight now available on STN
NEWS 19 Jan 29 Simultaneous left and right truncation added to COMPENDEX,
ENERGY, INSPEC
NEWS 20 Feb 13 CANCERLIT is no longer being updated
NEWS 21 Feb 24 METADEX enhancements
NEWS 22 Feb 24 PCTGEN now available on STN
NEWS 23 Feb 24 TEMA now available on STN
NEWS 24 Feb 26 NTIS now allows simultaneous left and right truncation
NEWS 25 Feb 26 PCTFULL now contains images
NEWS 26 Mar 04 SDI PACKAGE for monthly delivery of multifile SDI results
NEWS 27 Mar 19 APOLLIT offering free connect time in April 2003
NEWS 28 Mar 20 EVENTLINE will be removed from STN
NEWS 29 Mar 24 PATDPAFULL now available on STN
NEWS 30 Mar 24 Additional information for trade-named substances without
structures available in REGISTRY
NEWS 31 Apr 11 Display formats in DGENE enhanced
NEWS 32 Apr 14 MEDLINE Reload
NEWS 33 Apr 17 Polymer searching in REGISTRY enhanced
NEWS 34 Apr 21 Indexing from 1947 to 1956 being added to records in CA/CAPLUS
NEWS 35 Apr 21 New current-awareness alert (SDI) frequency in
WPIDS/WPINDEX/WPIX
NEWS 36 Apr 28 RDISCLOSURE now available on STN
NEWS 37 May 05 Pharmacokinetic information and systematic chemical names
added to PHAR

NEWS EXPRESS April 4 CURRENT WINDOWS VERSION IS V6.01a, CURRENT
MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),
AND CURRENT DISCOVER FILE IS DATED 01 APRIL 2003

NEWS HOURS STN Operating Hours Plus Help Desk Availability

NEWS INTER General Internet Information

NEWS LOGIN Welcome Banner and News Items

NEWS PHONE Direct Dial and Telecommunication Network Access to STN

NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 09:38:00 ON 11 MAY 2003

FILE 'REGISTRY' ENTERED AT 09:38:12 ON 11 MAY 2003
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2003

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=> s pyridopyrimidine/cn
L1 0 PYRIDOPYRIMIDINE/CN

=> s pyridopyrimidine

L2 O PYRIDOPYRIMIDINE

| | | |
|----------------------|------------|---------|
| => file caplus | SINCE FILE | TOTAL |
| COST IN U.S. DOLLARS | ENTRY | SESSION |
| FULL ESTIMATED COST | 9.64 | 9.85 |

FILE 'CAPLUS' ENTERED AT 09:39:49 ON 11 MAY 2003
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
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FILE COVERS 1907 - 11 May 2003 VOL 138 ISS 20
FILE LAST UPDATED: 9 May 2003 (20030509/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s wo200130768?/pn
L3 1 WO200130768?/PN
(WO2001030768/BN)

=> s wo-200116114?/pn
L4 1 WO-200116114?/PN
(WO2001016114/BN)

| | | |
|----------------------|------------|---------|
| => file reg | SINCE FILE | TOTAL |
| COST IN U.S. DOLLARS | ENTRY | SESSION |
| FULL ESTIMATED COST | 3.86 | 13.71 |

FILE 'REGISTRY' ENTERED AT 09:40:42 ON 11 MAY 2003
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 9 MAY 2003 HIGHEST RN 513416-44-9
DICTIONARY FILE UPDATES: 9 MAY 2003 HIGHEST RN 513416-44-9

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2003

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

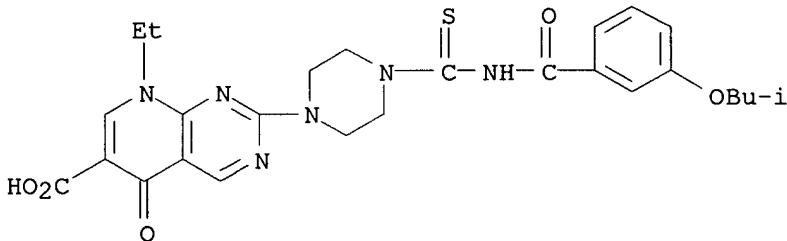
Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

```
=> s pyrido[2,3-d]pyrimidine/cn
L5          0 PYRIDO[2,3-D]PYRIMIDINE/CN

=> s pyrido[2,3-d]pyrimidine/cns
     140299 PYRIDO/CNS
     112135 2,3-D/CNS
     159578 PYRIMIDINE/CNS
L6          8193 PYRIDO[2,3-D]PYRIMIDINE/CNS
             ((PYRIDO(W)2,3-D(W)PYRIMIDINE)/CNS)

=> d scan cn str rid
'RID' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

L6  8193 ANSWERS  REGISTRY  COPYRIGHT 2003 ACS
IN  Pyrido[2,3-d]pyrimidine-6-carboxylic acid, 8-ethyl-5,8-dihydro-2-[4-
     [[[3-(2-methylpropoxy)benzoyl]amino]thioxomethyl]-1-piperazinyl]-5-oxo-
     (9CI)
MF  C26 H30 N6 O5 S
```



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

| | |
|--------|---|
| REG | - RN |
| SAM | - Index Name, MF, and structure - no RN |
| FIDE | - All substance data, except sequence data |
| IDE | - FIDE, but only 50 names |
| SQIDE | - IDE, plus sequence data |
| SQIDE3 | - Same as SQIDE, but 3-letter amino acid codes are used |
| SQD | - Protein sequence data, includes RN |
| SQD3 | - Same as SQD, but 3-letter amino acid codes are used |

SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties

EPROP - Table of experimental properties

PROP - EPROP and CALC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations

SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.

HELP FORMATS -- To see detailed descriptions of the predefined formats.

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> d scan cn str rsd

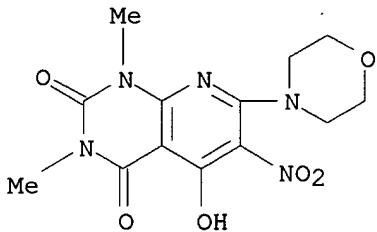
INVALID SCAN FIELD FOR FILE 'REGISTRY'

One or more of the display fields specified are not valid with DISPLAY SCAN in the current file. Enter HELP DSCAN at the arrow prompt (=>) for the list of fields that may be used when scanning the answers.

=> d cn str rsd 1 44 79

L6 ANSWER 1 OF 8193 REGISTRY COPYRIGHT 2003 ACS

CN Pyrido[2,3-d]pyrimidine-2,4(1H,3H)-dione, 5-hydroxy-1,3-dimethyl-7-(4-morpholinyl)-6-nitro- (9CI) (CA INDEX NAME)

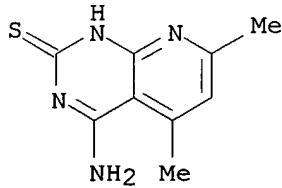


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

Ring System Data

| Elemental Analysis | Elemental Sequence | Size of the Rings | Ring System Formula | Identifier | Ring RID | Occurrence Count |
|--------------------|--------------------|-------------------|---------------------|------------|----------|------------------|
| EA | ES | SZ | RF | RID | | |
| C4NO | NC2OC2 | 6 | C4NO | 46.402.1 | 1 | |
| C4N2-C5N | NCNC3-NC5 | 6-6 | C7N3 | 591.304.18 | 1 | |

L6 ANSWER 44 OF 8193 REGISTRY COPYRIGHT 2003 ACS
CN Pyrido[2,3-d]pyrimidine-2(1H)-thione, 4-amino-5,7-dimethyl- (9CI)
(CA INDEX NAME)



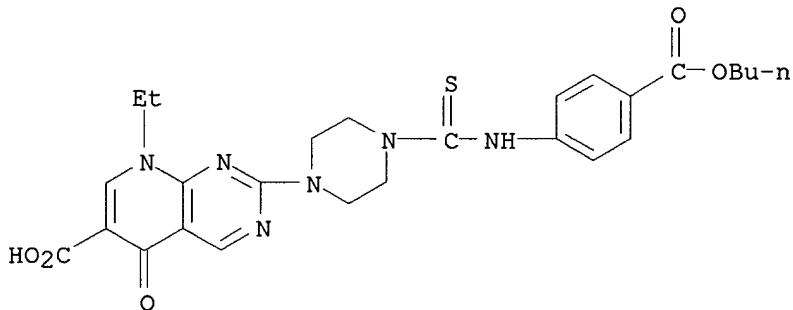
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

Ring System Data

| Elemental Analysis | Elemental Sequence | Size of the Rings | Ring System Formula | Identifier | Ring RID | Occurrence Count |
|--------------------|--------------------|-------------------|---------------------|------------|----------|------------------|
| EA | ES | SZ | RF | RID | RID | Count |
| C4N2-C5N | NCNC3-NC5 6-6 | C7N3 | 1591.304.27 1 | | | |

L6 ANSWER 79 OF 8193 REGISTRY COPYRIGHT 2003 ACS
CN Pyrido[2,3-d]pyrimidine-6-carboxylic acid, 2-[4-[[[4-

(butoxycarbonyl)phenyl]amino]thioxomethyl]-1-piperazinyl]-8-ethyl-5,8-dihydro-5-oxo- (9CI) (CA INDEX NAME)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

Ring System Data

| Elemental Analysis | Elemental Sequence | Size of the Rings | Ring System Formula | Identifier | Occurrence RID | Count |
|--------------------|--------------------|-------------------|---------------------|------------|----------------|-------|
| EA | ES | SZ | RF | RID | | |
| C6 | C6 | 6 | C6 | 46.150.18 | 1 | |
| C4N2 | NC2NC2 | 6 | C4N2 | 46.383.1 | 1 | |
| C4N2-C5N | NCNC3-NC5 | 6-6 | C7N3 | 591.304.13 | 1 | |

=> s 13 and 591.304/rid
0 WO200130768?/PN
(WO2001030768/BN)
14204 591.304/RID
L7 0 L3 AND 591.304/RID

=> transfer 13
ENTER ANSWER NUMBERS, RANGES (1-), OR ?:.
ENTER DISPLAY FIELDS (TI) OR ?:rn
L8 TRANSFER L3 1- RN : 427 TERMS
L9 427 L8

=> s 19 and 591.304/rid
14204 591.304/RID
L10 0 L9 AND 591.304/RID

=> transfer 14 1- rn
L11 TRANSFER L4 1- RN : 46 TERMS
L12 46 L11

=> s 112 and 591.304/rid
14204 591.304/RID

L13 0 L12 AND 591.304/RID

=> s l12 and 591/rid
1877036 591/RID

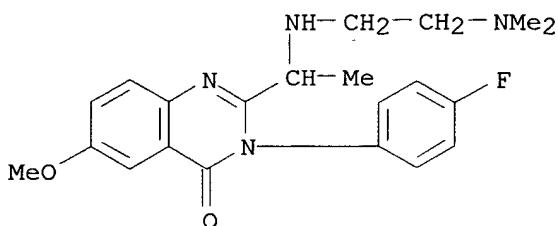
L14 25 L12 AND 591/RID

=> s l9 and 591/rid
1877036 591/RID

L15 420 L9 AND 591/RID

=> d l14 cn str cbib pi

L14 ANSWER 1 OF 25 REGISTRY COPYRIGHT 2003 ACS
CN 4(3H)-Quinazolinone, 2-[1-[[2-(dimethylamino)ethyl]amino]ethyl]-3-(4-fluorophenyl)-6-methoxy- (9CI) (CA INDEX NAME)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

REFERENCE 1: 134:222723 Preparation of quinazolinones for modulating CXR3 function. Schall, Thomas J.; Dairaghi, Daniel J.; McMaster, Brian E. (Chemocentryx, Inc., USA). PCT Int. Appl. WO 2001016114 A2 20010308, 53 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 2000-US23556 20000825. PRIORITY: US 1999-PV151212 19990827.

| | PATENT NO. | KIND | DATE | APPLICATION NO. | DATE | |
|----|---------------|---|----------|-----------------|----------|--|
| PI | WO 2001016114 | A2 | 20010308 | WO 2000-US23556 | 20000825 | |
| | W: | AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. | | | | |
| | EP 1216232 | A1 | 20020626 | EP 2000-959489 | 20000825 | |
| | R: | AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL | | | | |

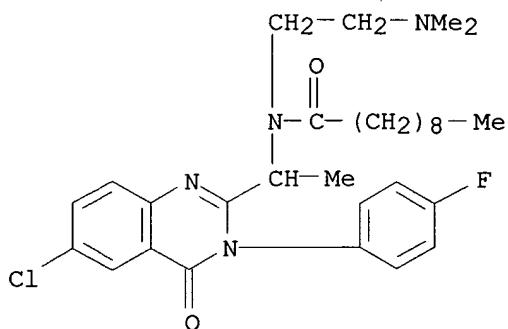
US 6559160

B1 20030506

US 2000-648329 20000825

=> d scan 114

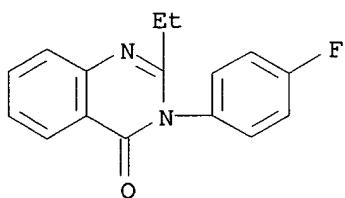
L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[1-[6-chloro-3-(4-fluorophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]ethyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C30 H40 Cl F N4 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

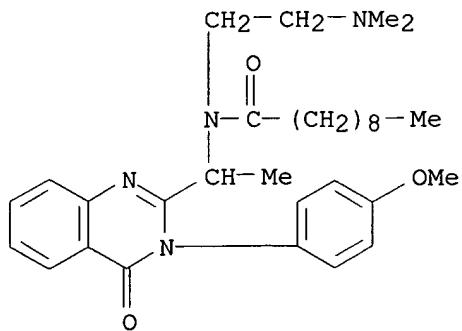
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):24

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-ethyl-3-(4-fluorophenyl)- (9CI)
MF C16 H13 F N2 O



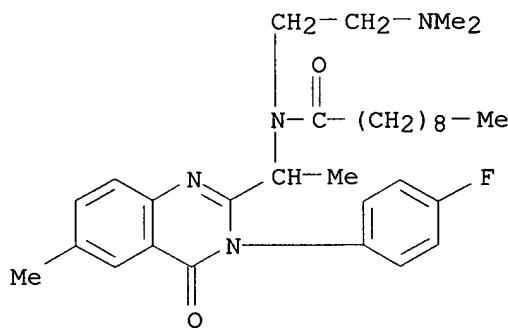
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[1-[3,4-dihydro-3-(4-methoxyphenyl)-4-oxo-2-quinazolinyl]ethyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C31 H44 N4 O3



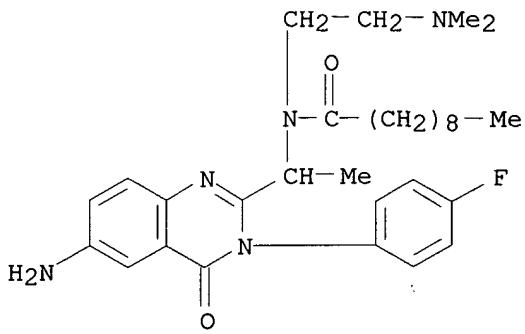
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-6-methyl-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C31 H43 F N4 O2



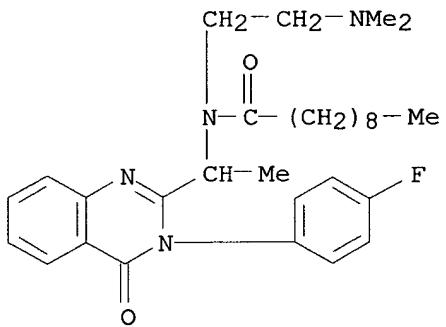
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[1-[6-amino-3-(4-fluorophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]ethyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C30 H42 F N5 O2



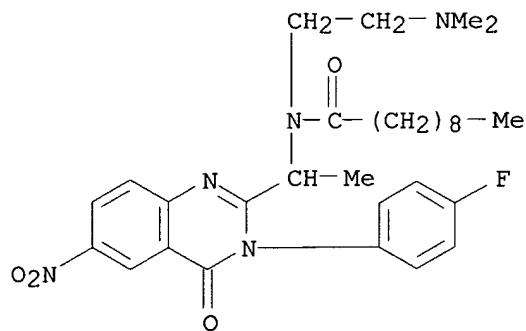
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C30 H41 F N4 O2



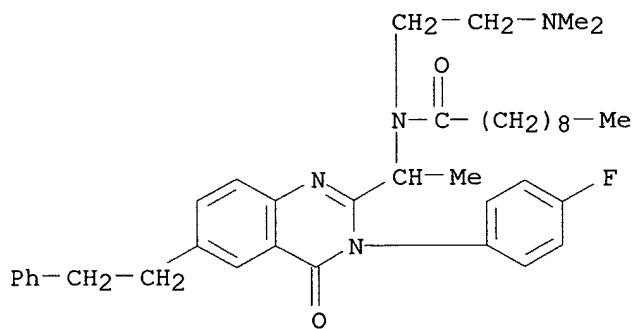
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-6-nitro-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C30 H40 F N5 O4



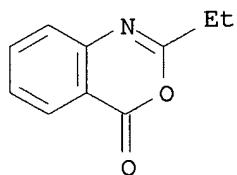
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxo-6-(2-phenylethyl)-2-quinazolinyl]ethyl]- (9CI)
MF C38 H49 F N4 O2



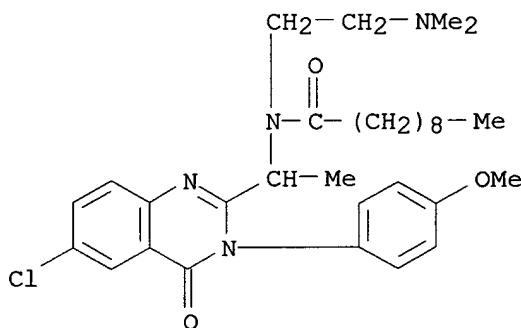
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4H-3,1-Benzoxazin-4-one, 2-ethyl- (7CI, 8CI, 9CI)
MF C10 H9 N O2



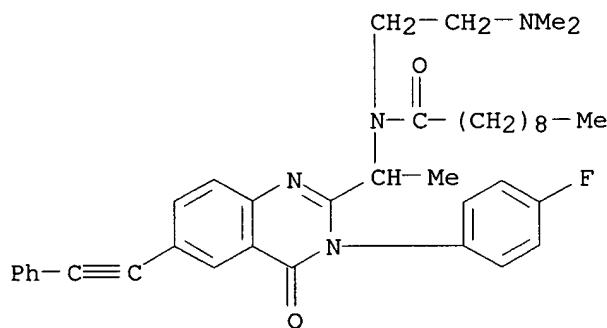
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[1-[6-chloro-3,4-dihydro-3-(4-methoxyphenyl)-4-oxo-2-quinazolinyl]ethyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C31 H43 Cl N4 O3



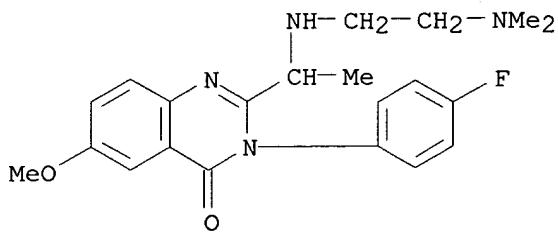
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxo-6-(phenylethyynyl)-2-quinazolinyl]ethyl]- (9CI)
MF C38 H45 F N4 O2



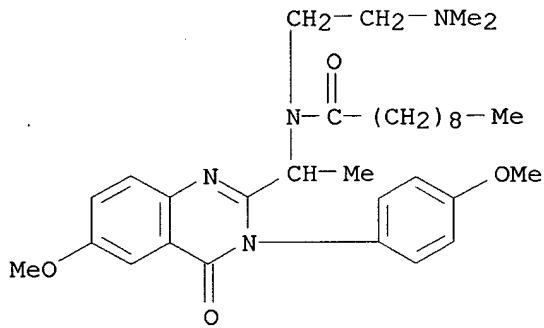
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-[1-[(2-(dimethylamino)ethyl)amino]ethyl]-3-(4-fluorophenyl)-6-methoxy- (9CI)
MF C21 H25 F N4 O2



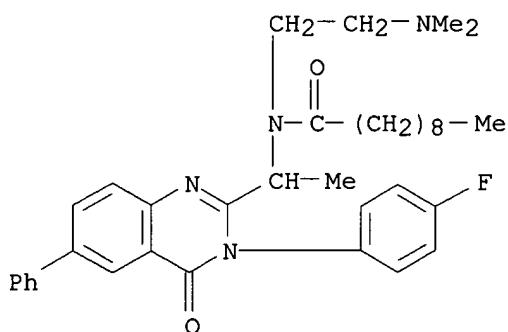
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[1-[3,4-dihydro-6-methoxy-3-(4-methoxyphenyl)-4-oxo-2-quinazolinyl]ethyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C32 H46 N4 O4



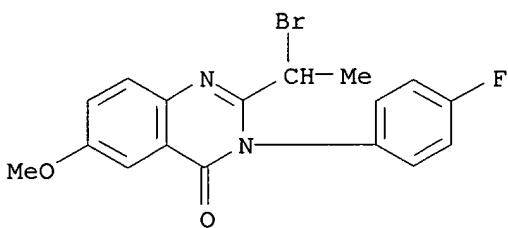
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxo-6-phenyl-2-quinazolinyl]ethyl]- (9CI)
MF C36 H45 F N4 O2



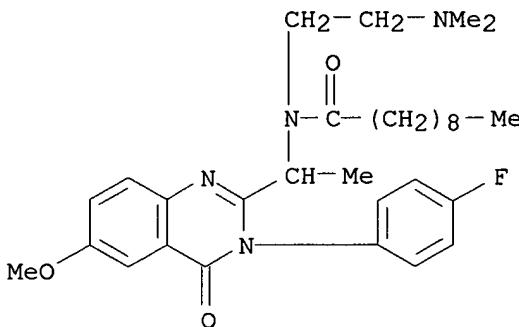
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-(1-bromoethyl)-3-(4-fluorophenyl)-6-methoxy- (9CI)
MF C17 H14 Br F N2 O2



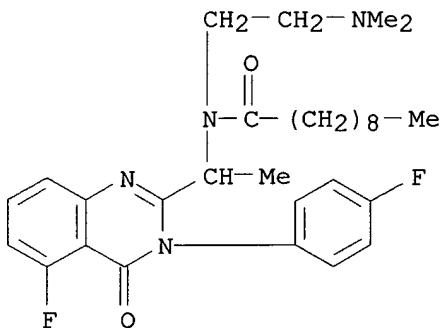
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-6-methoxy-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C31 H43 F N4 O3



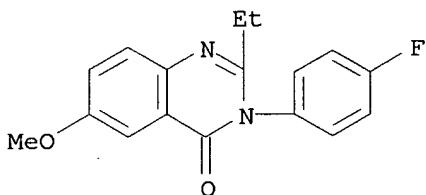
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[5-fluoro-3-(4-fluorophenyl)-
3,4-dihydro-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C30 H40 F2 N4 O2



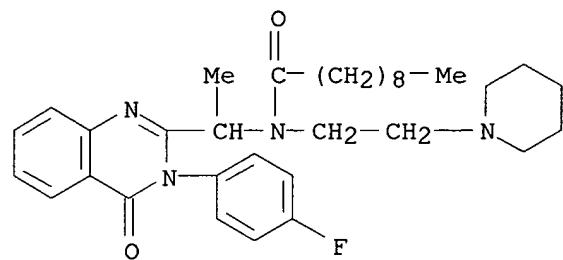
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-ethyl-3-(4-fluorophenyl)-6-methoxy- (9CI)
MF C17 H15 F N2 O2



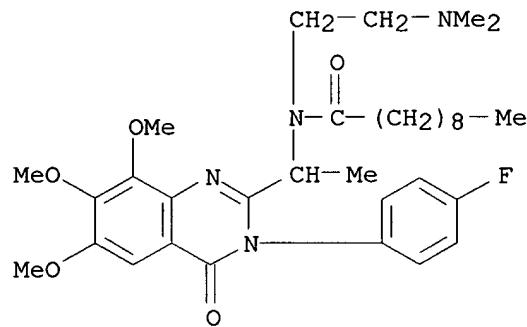
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxo-2-
quinazolinyl]ethyl]-N-[2-(1-piperidinyl)ethyl]- (9CI)
MF C33 H45 F N4 O2



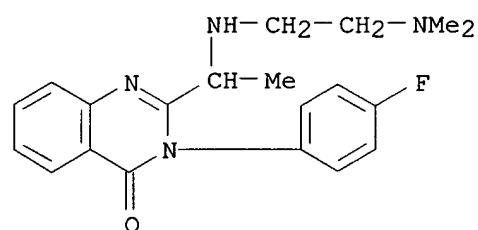
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-6,7,8-trimethoxy-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C33 H47 F N4 O5



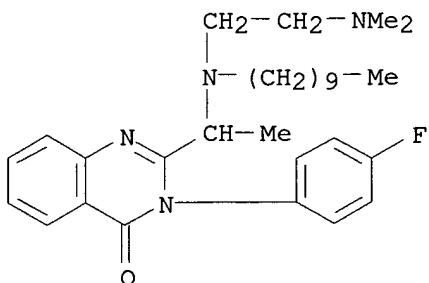
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-[1-[[2-(dimethylamino)ethyl]amino]ethyl]-3-(4-fluorophenyl)- (9CI)
MF C20 H23 F N4 O



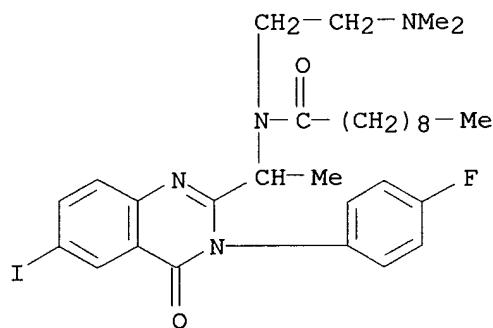
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-[1-[decyl[2-(dimethylamino)ethyl]amino]ethyl]-3-(4-fluorophenyl)- (9CI)
MF C30 H43 F N4 O



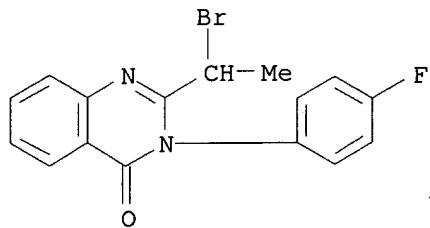
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[2-(dimethylamino)ethyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-6-iodo-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C30 H40 F I N4 O2



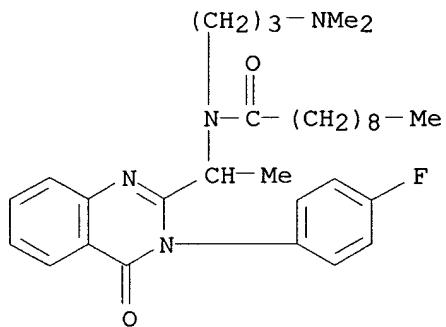
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-(1-bromoethyl)-3-(4-fluorophenyl)- (9CI)
MF C16 H12 Br F N2 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 25 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Decanamide, N-[3-(dimethylamino)propyl]-N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]ethyl]- (9CI)
MF C31 H43 F N4 O2

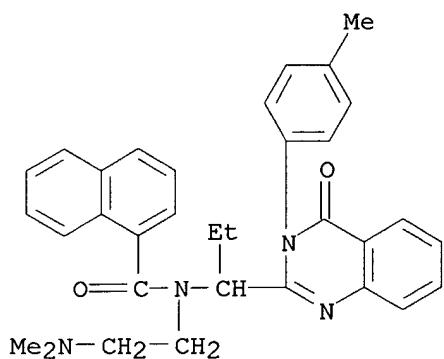


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> d 115 cn str cbib pi

L15 ANSWER 1 OF 420 REGISTRY COPYRIGHT 2003 ACS
CN 1-Naphthalenecarboxamide, N-[1-[3,4-dihydro-3-(4-methylphenyl)-4-oxo-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]- (9CI) (CA INDEX NAME)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

REFERENCE 1: 136:53759 Preparation of N-acylquinazolinonealkylamines as KSP kinesin inhibitors. Finer, Jeffrey T.; Bergnes, Gustav; Feng, Bainian; Smith, Whitney W.; Chabala, John C.; Morgans, David J., Jr. (Cytokinetics, Inc., USA). PCT Int. Appl. WO 2001098278 A1 20011227, 179 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR. (English). CODEN: PIXXD2. APPLICATION: WO 2001-US13901 20010427. PRIORITY: US 2000-PV213104 20000621; US 2000-699047 20001024.

| PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|---------------|--|-------------|-----------------|----------|
| PI | WO 2001098278 | A1 20011227 | WO 2001-US13901 | 20010427 |
| | W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR. | | | |
| US 6545004 | B1 20030408 | | US 2000-699047 | 20001024 |
| JP 2003048881 | A2 20030221 | | JP 2002-156766 | 20001026 |
| EP 1296959 | A1 20030402 | | EP 2001-932769 | 20010427 |
| NO 2002006172 | A 20030220 | | NO 2002-6172 | 20021220 |

REFERENCE 2: 134:326543 Methods and compositions utilizing quinazolinones as KSP kinesin modulators. Finer, Jeffrey T.; Bergnes, Gustave; Feng, Bainian; Smith, Whitney W.; Chabala, John C. (Cytokinetics, Inc., USA). PCT Int. Appl. WO 2001030768 A1 20010503, 168 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK,

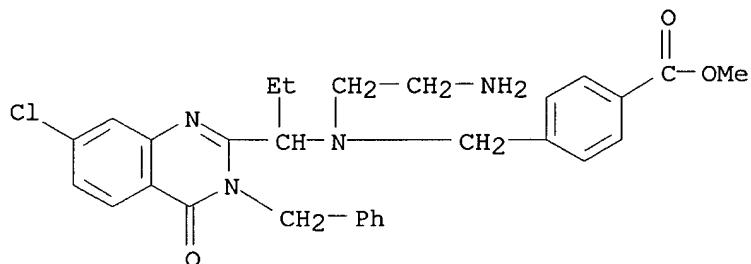
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 GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG.
 (English). CODEN: PIXXD2. APPLICATION: WO 2000-US29585 20001026.

PRIORITY: US 1999-PV198253 19991027; US 2000-PV213104 20000621.

| PI | PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|----|---------------|---|----------|-----------------|----------|
| | WO 2001030768 | A1 | 20010503 | WO 2000-US29585 | 20001026 |
| | W: | AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
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SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN,
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CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG | | | |
| | BR 2000015110 | A | 20020702 | BR 2000-15110 | 20001026 |
| | EP 1226129 | A1 | 20020731 | EP 2000-976656 | 20001026 |
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IE, SI, LT, LV, FI, RO, MK, CY, AL | | | |
| | JP 2003048881 | A2 | 20030221 | JP 2002-156766 | 20001026 |
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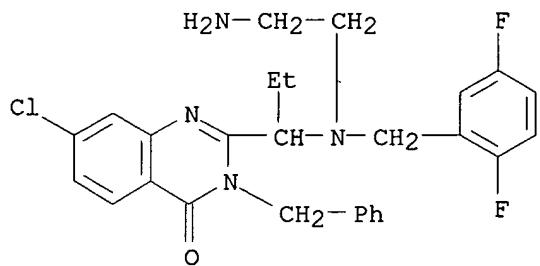
L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Benzoic acid, 4-[(2-aminoethyl)[1-[7-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]amino]methyl]-, methyl ester (9CI)
 MF C29 H31 Cl N4 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):25

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 4(3H)-Quinazolinone, 2-[1-[(2-aminoethyl)[(2,5-difluorophenyl)methyl]amino]propyl]-7-chloro-3-(phenylmethyl)- (9CI)
 MF C27 H27 Cl F2 N4 O

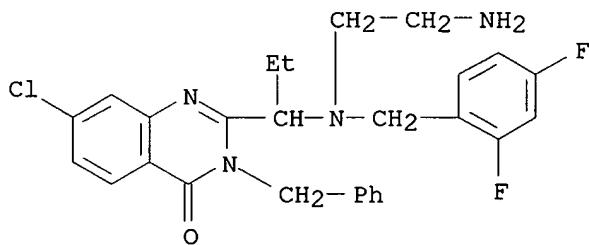


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS

IN 4(3H)-Quinazolinone, 2-[1-[(2-aminoethyl)[(2,4-difluorophenyl)methyl]amino]propyl]-7-chloro-3-(phenylmethyl)-(9CI)

MF C27 H27 Cl F2 N4 O

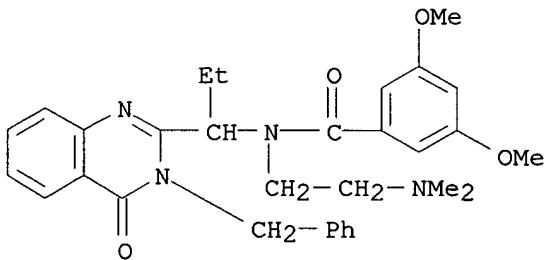


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L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS

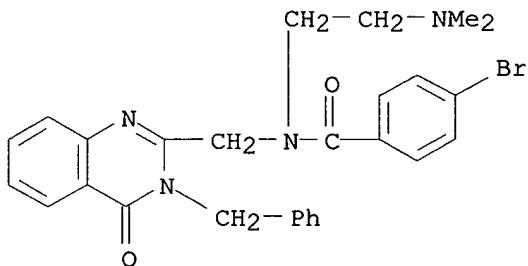
IN Benzamide, N-[1-[3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]-3,5-dimethoxy-(9CI)

MF C31 H36 N4 O4



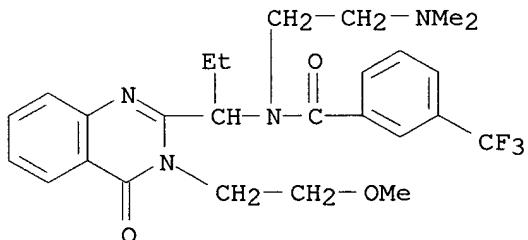
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, 4-bromo-N-[[3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]methyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C27 H27 Br N4 O2



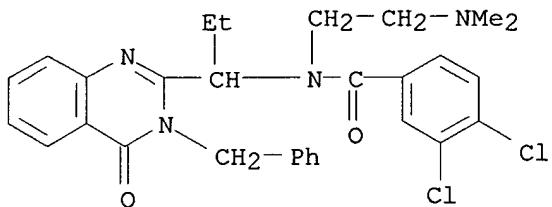
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[1-[3,4-dihydro-3-(2-methoxyethyl)-4-oxo-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]-3-(trifluoromethyl)- (9CI)
MF C26 H31 F3 N4 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

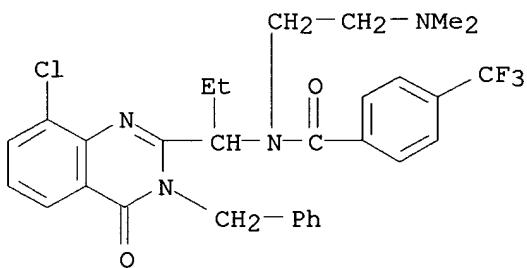
L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, 3,4-dichloro-N-[1-[3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C29 H30 Cl2 N4 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS

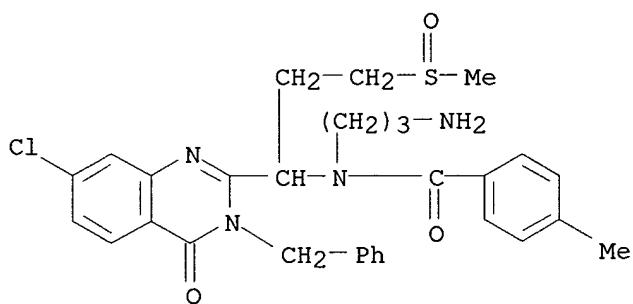
IN Benzamide, N-[1-[8-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]-4-(trifluoromethyl)- (9CI)
MF C30 H30 Cl F3 N4 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

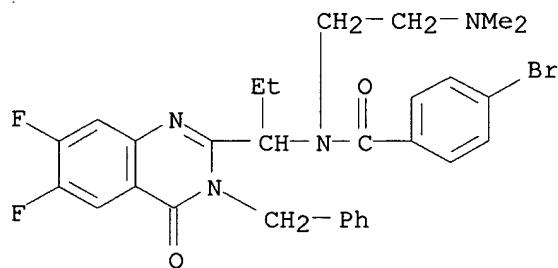
L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS

IN Benzamide, N-(3-aminopropyl)-N-[1-[7-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]-3-(methylsulfinyl)propyl]-4-methyl- (9CI)
MF C30 H33 Cl N4 O3 S



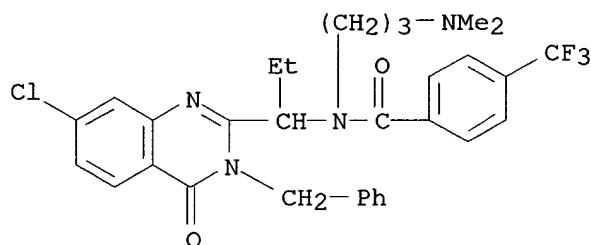
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, 4-bromo-N-[1-[6,7-difluoro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C29 H29 Br F2 N4 O2



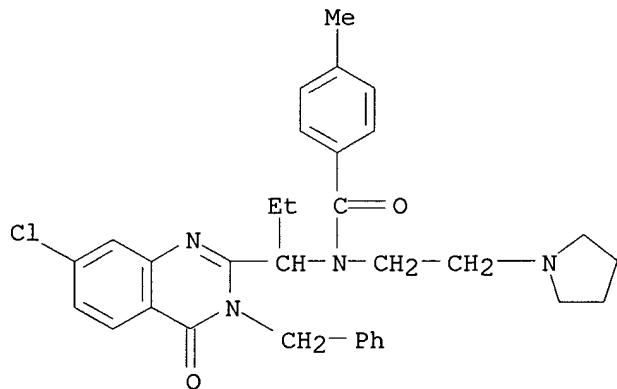
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[1-[7-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-N-[3-(dimethylamino)propyl]-4-(trifluoromethyl)- (9CI)
MF C31 H32 Cl F3 N4 O2



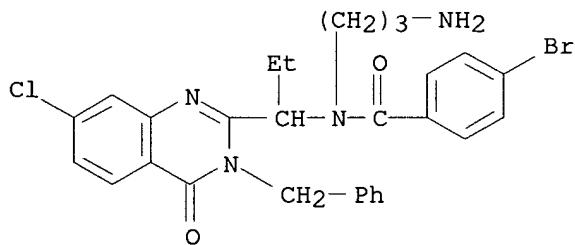
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[1-[7-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-4-methyl-N-[2-(1-pyrrolidinyl)ethyl]- (9CI)
MF C32 H35 Cl N4 O2



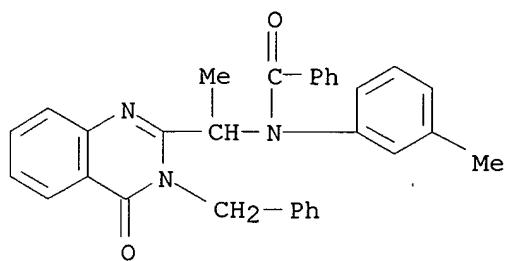
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-(3-aminopropyl)-4-bromo-N-[1-[7-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]- (9CI)
MF C28 H28 Br Cl N4 O2



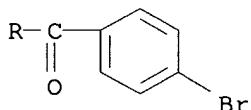
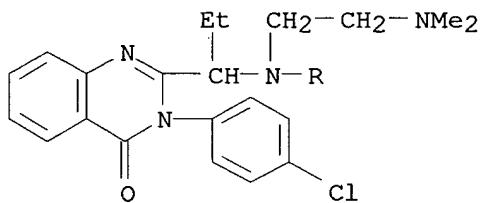
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[1-[3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]ethyl]-N-(3-methylphenyl)- (9CI)
MF C31 H27 N3 O2



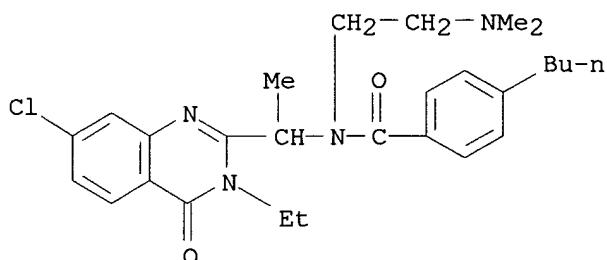
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, 4-bromo-N-[1-[3-(4-chlorophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C28 H28 Br Cl N4 O2



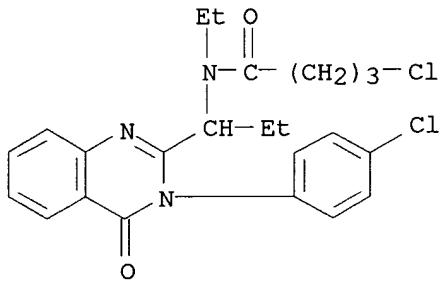
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, 4-butyl-N-[1-(7-chloro-3-ethyl-3,4-dihydro-4-oxo-2-quinazolinyl)ethyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C27 H35 Cl N4 O2



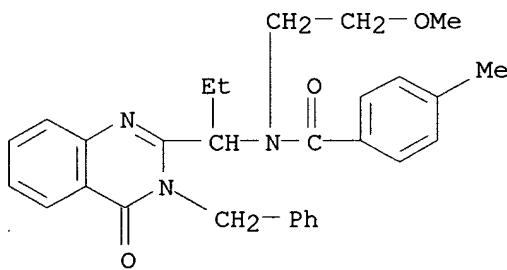
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Butanamide, 4-chloro-N-[1-[3-(4-chlorophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]propyl]-N-ethyl- (9CI)
MF C23 H25 Cl2 N3 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

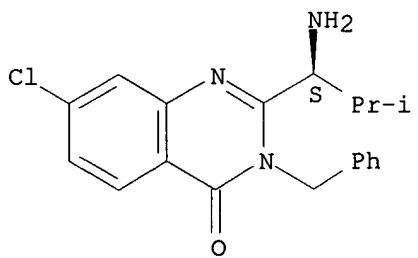
L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[1-[3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]-N-(2-methoxyethyl)-4-methyl- (9CI)
MF C29 H31 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

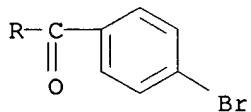
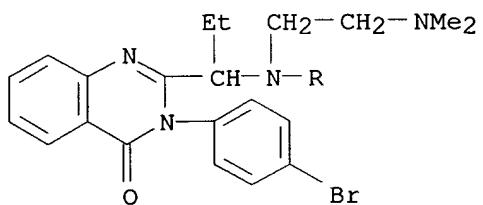
L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-[(1S)-1-amino-2-methylpropyl]-7-chloro-3-(phenylmethyl)- (9CI)
MF C19 H20 Cl N3 O

Absolute stereochemistry.



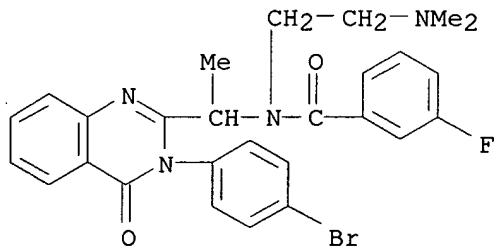
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, 4-bromo-N-[1-[3-(4-bromophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]propyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C28 H28 Br2 N4 O2



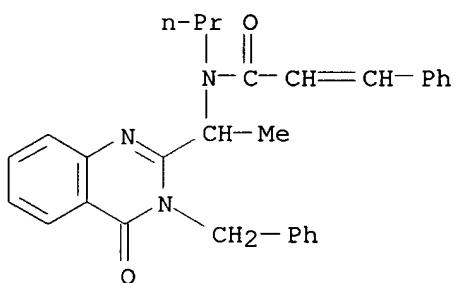
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[1-[3-(4-bromophenyl)-3,4-dihydro-4-oxo-2-quinazolinyl]ethyl]-N-[2-(dimethylamino)ethyl]-3-fluoro- (9CI)
MF C27 H26 Br F N4 O2



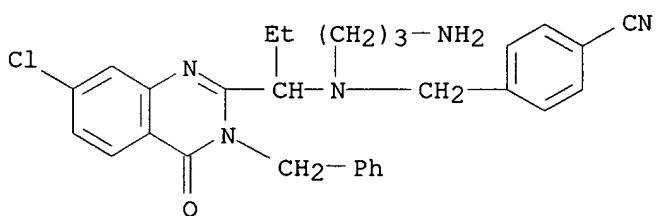
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 2-Propenamide, N-[1-[3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]ethyl]-3-phenyl-N-propyl- (9CI)
MF C29 H29 N3 O2



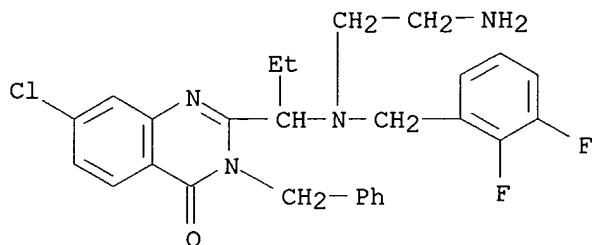
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzonitrile, 4-[(3-aminopropyl)[1-[7-chloro-3,4-dihydro-4-oxo-3-(phenylmethyl)-2-quinazolinyl]propyl]amino]methyl- (9CI)
MF C29 H30 Cl N5 O



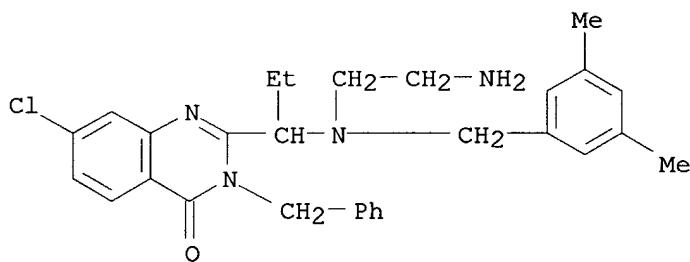
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-[1-[(2-aminoethyl)[(2,3-difluorophenyl)methyl]amino]propyl]-7-chloro-3-(phenylmethyl)- (9CI)
MF C27 H27 Cl F2 N4 O



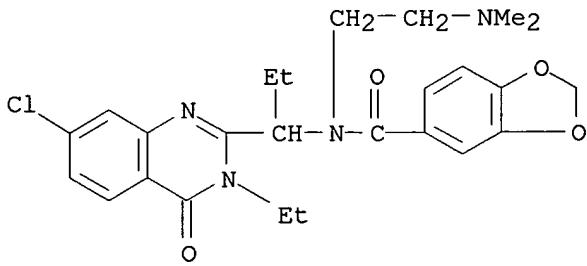
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4(3H)-Quinazolinone, 2-[1-[(2-aminoethyl)[(3,5-dimethylphenyl)methyl]amino]propyl]-7-chloro-3-(phenylmethyl)- (9CI)
MF C29 H33 Cl N4 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L15 420 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 1,3-Benzodioxole-5-carboxamide, N-[1-(7-chloro-3-ethyl-3,4-dihydro-4-oxo-2-quinazolinyl)propyl]-N-[2-(dimethylamino)ethyl]- (9CI)
MF C25 H29 Cl N4 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> logoff
ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF
LOGOFF? (Y)/N/HOLD:.

| COST IN U.S. DOLLARS | SINCE FILE
ENTRY | TOTAL
SESSION |
|----------------------|---------------------|------------------|
| FULL ESTIMATED COST | 21.94 | 95.55 |

STN INTERNATIONAL LOGOFF AT 09:52:54 ON 11 MAY 2003

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: ssspta1611txm

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * * * Welcome to STN International * * * * * * * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 Apr 08 "Ask CAS" for self-help around the clock
NEWS 3 Jun 03 New e-mail delivery for search results now available
NEWS 4 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN
NEWS 5 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE)
now available on STN
NEWS 6 Aug 26 Sequence searching in REGISTRY enhanced
NEWS 7 Sep 03 JAPIO has been reloaded and enhanced
NEWS 8 Sep 16 Experimental properties added to the REGISTRY file
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NEWS 10 Oct 01 CASREACT Enriched with Reactions from 1907 to 1985
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NEWS 14 Nov 25 More calculated properties added to REGISTRY
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NEWS 16 Dec 17 PCTFULL now covers WP/PCT Applications from 1978 to date
NEWS 17 Dec 17 TOXCENTER enhanced with additional content
NEWS 18 Dec 17 Adis Clinical Trials Insight now available on STN
NEWS 19 Jan 29 Simultaneous left and right truncation added to COMPENDEX,
ENERGY, INSPEC
NEWS 20 Feb 13 CANCERLIT is no longer being updated
NEWS 21 Feb 24 METADEX enhancements
NEWS 22 Feb 24 PCTGEN now available on STN
NEWS 23 Feb 24 TEMA now available on STN
NEWS 24 Feb 26 NTIS now allows simultaneous left and right truncation
NEWS 25 Feb 26 PCTFULL now contains images
NEWS 26 Mar 04 SDI PACKAGE for monthly delivery of multifile SDI results
NEWS 27 Mar 19 APOLLIT offering free connect time in April 2003
NEWS 28 Mar 20 EVENTLINE will be removed from STN
NEWS 29 Mar 24 PATDPAFULL now available on STN
NEWS 30 Mar 24 Additional information for trade-named substances without
structures available in REGISTRY
NEWS 31 Apr 11 Display formats in DGENE enhanced
NEWS 32 Apr 14 MEDLINE Reload
NEWS 33 Apr 17 Polymer searching in REGISTRY enhanced
NEWS 34 Apr 21 Indexing from 1947 to 1956 being added to records in CA/CAPLUS
NEWS 35 Apr 21 New current-awareness alert (SDI) frequency in
WPIDS/WPINDEX/WPIX
NEWS 36 Apr 28 RDISCLOSURE now available on STN
NEWS 37 May 05 Pharmacokinetic information and systematic chemical names
added to PHAR

NEWS EXPRESS April 4 CURRENT WINDOWS VERSION IS V6.01a, CURRENT
MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),
AND CURRENT DISCOVER FILE IS DATED 01 APRIL 2003

NEWS HOURS STN Operating Hours Plus Help Desk Availability
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NEWS WWW CAS World Wide Web Site (general information)

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Thomas McKenzie 05/11/2003

FULL ESTIMATED COST

0.21

0.21

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provided by InfoChem.

STRUCTURE FILE UPDATES: 9 MAY 2003 HIGHEST RN 513416-44-9
DICTIONARY FILE UPDATES: 9 MAY 2003 HIGHEST RN 513416-44-9

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2003

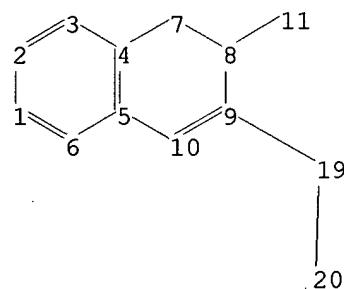
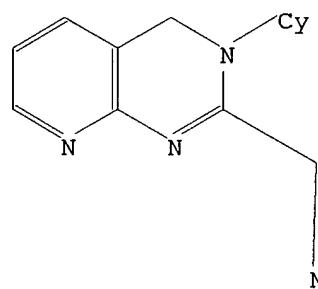
Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>.

=>

Uploading C:\Program Files\Stnexp\Queries\10015532.str



chain nodes :
11 12 13 14 19 20
ring nodes :

<10/015, 532Page 35

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1 2 3 4 5 6 7 8 9 10
chain bonds :
8-11 9-19 13-14 19-20
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 9-10
exact/norm bonds :
4-7 5-10 7-8 8-9 8-11 9-10 13-14 19-20
exact bonds :
9-19
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
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Match level :

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1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:CLASS 13:CLASS 14:CLASS 19:CLASS 20:CLASS
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L1 STRUCTURE UPLOADED

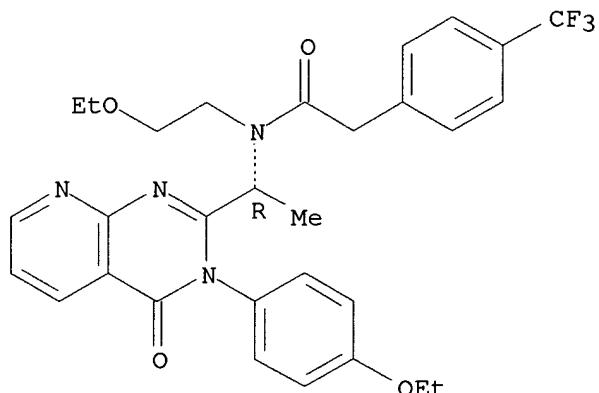
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SAMPLE SEARCH INITIATED 10:08:03 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 92 TO ITERATE
100.0% PROCESSED 92 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
                      BATCH **COMPLETE**
PROJECTED ITERATIONS: 1265 TO 2415
PROJECTED ANSWERS: 1 TO 80
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L2 1 SEA SSS SAM L1

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=> d scan
L2 1 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzeneacetamide, N-(2-ethoxyethyl)-N-[(1R)-1-[3-(4-ethoxyphenyl)-3,4-
dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-(trifluoromethyl)- (9CI)
MF C30 H31 F3 N4 O4
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Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

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FULL SEARCH INITIATED 10:08:24 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2262 TO ITERATE

100.0% PROCESSED 2262 ITERATIONS 21 ANSWERS
SEARCH TIME: 00.00.01

L3 21 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS SINCE FILE TOTAL
FULL ESTIMATED COST ENTRY SESSION
148.15 148.36

FILE 'CAPLUS' ENTERED AT 10:08:31 ON 11 MAY 2003
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FILE COVERS 1907 - 11 May 2003 VOL 138 ISS 20
FILE LAST UPDATED: 9 May 2003 (20030509/ED)

This file contains CAS Registry Numbers for easy and accurate

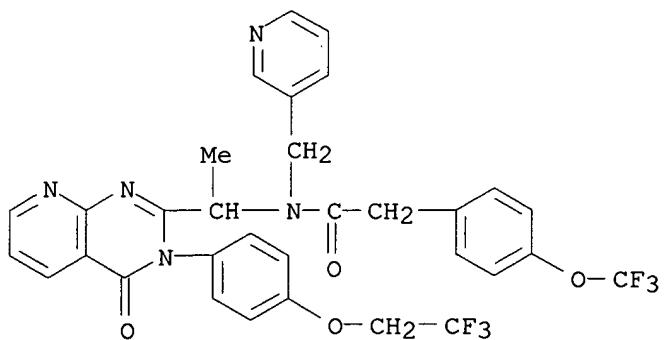
substance identification.

=> s 13
L4 2 L3

=> d 1-2 cbib pi hitstr

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2003 ACS
2002:813938 Document No. 137:337907 Preparation of N-(heteroarylalkyl)acylamides as CXCR3 antagonists for treatment of inflammatory or immune conditions. Medina, Julio C.; Johnson, Michael G.; Li, An-Rong; Liu, Jiwen; Huang, Alan Xi; Zhu, Liusheng; Marcus, Andrew P. (Tularik Inc., USA). PCT Int. Appl. WO 2002083143 A1 20021024, 205 pp.
DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR. (English). CODEN: PIXXD2.
APPLICATION: WO 2001-US47850 20011211. PRIORITY: US 2000-PV255241 20001211; US 2001-PV296499 20010606.

| | PATENT NO. | KIND | DATE | APPLICATION NO. | DATE | | |
|----|---|--|---------------|-----------------|----------|---------------|----------|
| PI | WO 2002083143 | A1 | 20021024 | WO 2001-US47850 | 20011211 | | |
| | W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM | RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG | US 2002169159 | A1 | 20021114 | US 2001-15532 | 20011211 |
| | US 2003069234 | A1 | 20030410 | US 2002-164690 | 20020606 | | |
| | US 2003055054 | A1 | 20030320 | US 2002-231895 | 20020829 | | |
| IT | 473720-05-7P | | | | | | |
| | RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses) | (CXCR3 antagonist; preparation of N-(heteroarylalkyl)acylamides as CXCR3 antagonists for treatment of inflammatory or immune conditions) | | | | | |
| RN | 473720-05-7 | CAPLUS | | | | | |
| CN | Benzeneacetamide, N-[1-[3,4-dihydro-4-oxo-3-[4-(2,2,2-trifluoroethoxy)phenyl]pyrido[2,3-d]pyrimidin-2-yl]ethyl]-N-(3-pyridinylmethyl)-4-(trifluoromethoxy)-(9CI) (CA INDEX NAME) | | | | | | |



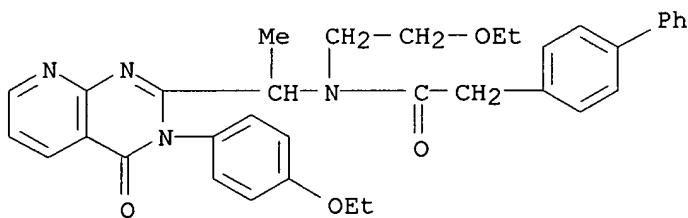
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473720-16-0P 473720-18-2P 473720-19-3P
473720-20-6P 473720-21-7P 473720-22-8P
473720-23-9P 473720-24-0P 473720-25-1P
473720-26-2P 473720-30-8P 473722-68-8P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(CXCR3 antagonist; preparation of N-(heteroarylalkyl)acylamides as CXCR3 antagonists for treatment of inflammatory or immune conditions)

RN 473718-93-3 CAPPLUS

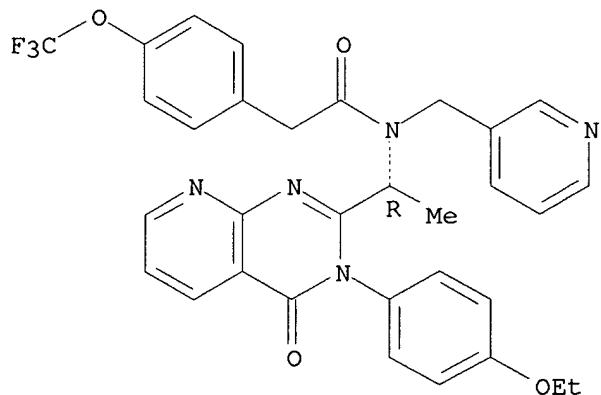
CN [1,1'-Biphenyl]-4-acetamide, N-(2-ethoxyethyl)-N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]- (9CI) (CA INDEX NAME)



RN 473719-41-4 CAPPLUS

CN Benzeneacetamide, N-[(1R)-1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-N-(3-pyridinylmethyl)-4-(trifluoromethoxy)- (9CI) (CA INDEX NAME)

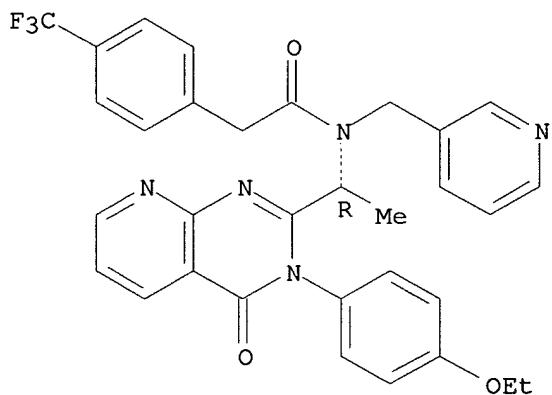
Absolute stereochemistry.



RN 473719-45-8 CAPLUS

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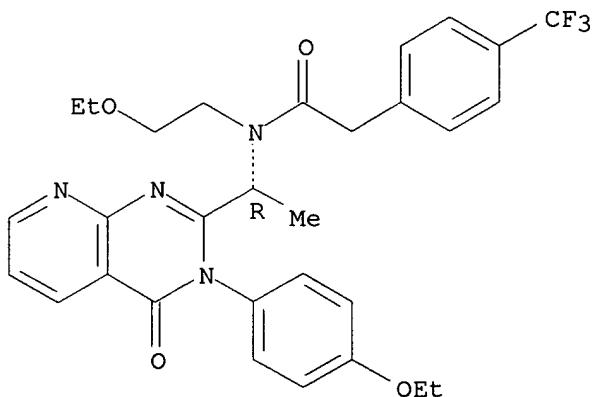
Absolute stereochemistry.



RN 473719-49-2 CAPLUS

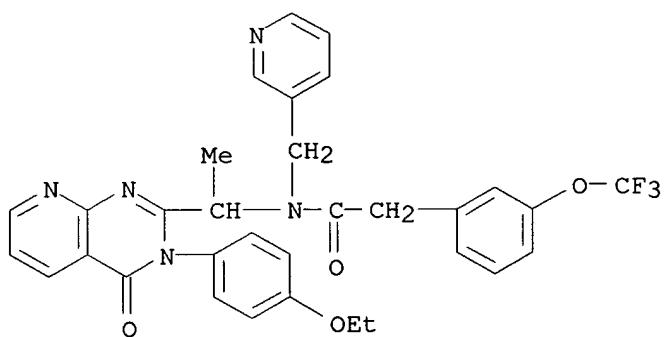
CN Benzeneacetamide, N-(2-ethoxyethyl)-N-[(1R)-1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-(trifluoromethyl)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



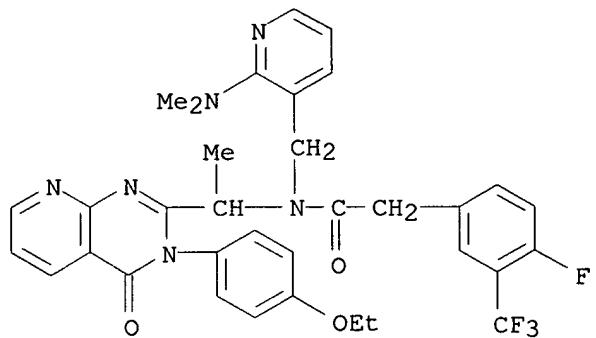
RN 473719-90-3 CAPLUS

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(CA INDEX NAME)



RN 473719-95-8 CAPLUS

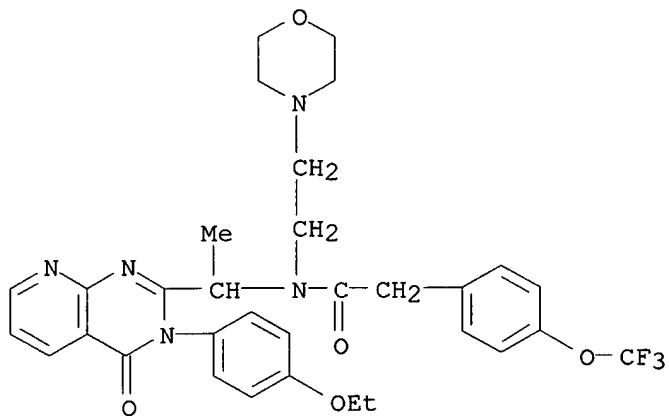
CN Benzeneacetamide, N-[2-(dimethylamino)-3-pyridinylmethyl]-N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-3-(trifluoromethyl)- (9CI) (CA INDEX NAME)



RN 473720-16-0 CAPLUS

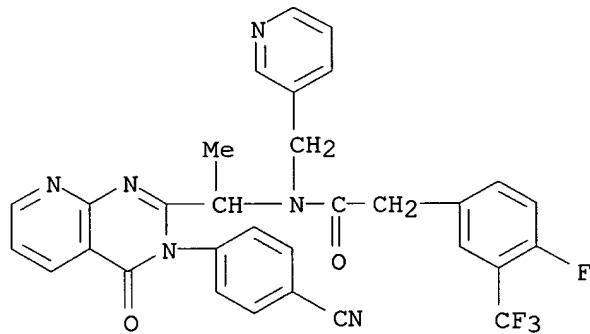
Thomas McKenzie 05/11/2003

CN Benzeneacetamide, N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-N-[2-(4-morpholinyl)ethyl]-4-(trifluoromethoxy)-(9CI) (CA INDEX NAME)



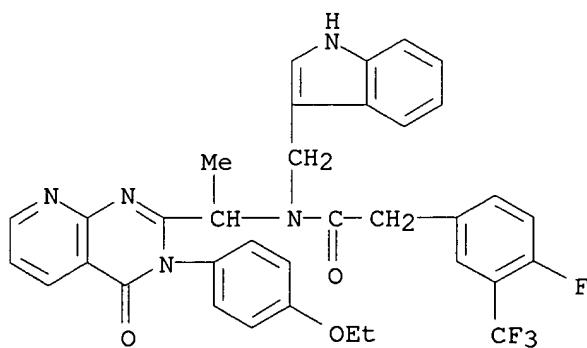
RN 473720-18-2 CAPLUS

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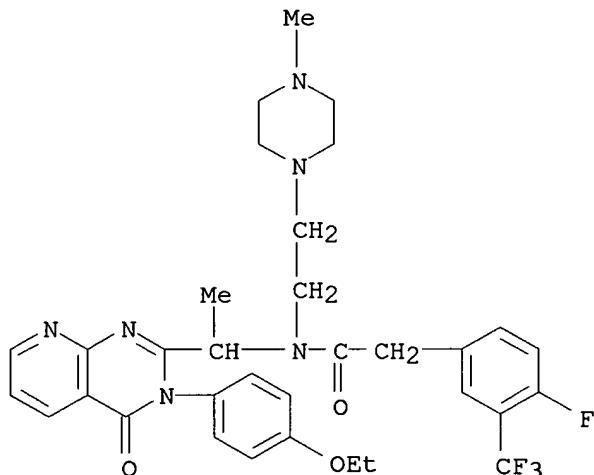
RN 473720-19-3 CAPLUS

CN Benzeneacetamide, N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-N-(1H-indol-3-ylmethyl)-3-(trifluoromethyl)-(9CI) (CA INDEX NAME)



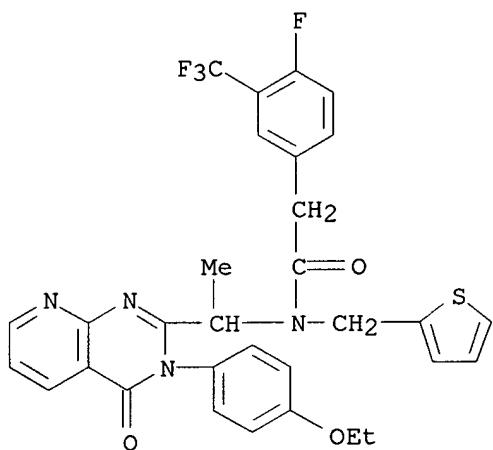
RN 473720-20-6 CAPLUS

CN Benzeneacetamide, N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-N-[2-(4-methyl-1-piperazinyl)ethyl]-3-(trifluoromethyl)- (9CI) (CA INDEX NAME)



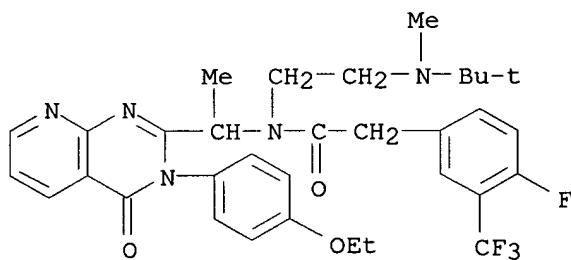
RN 473720-21-7 CAPLUS

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RN 473720-22-8 CAPLUS

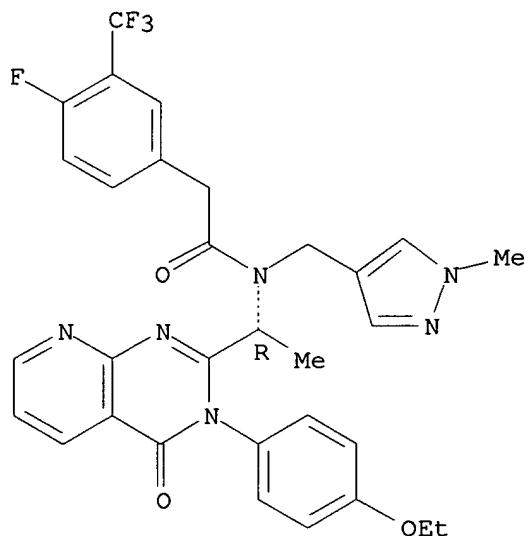
CN Benzeneacetamide, N-[2-[(1,1-dimethylethyl)methylamino]ethyl]-N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-3-(trifluoromethyl)- (9CI) (CA INDEX NAME)



RN 473720-23-9 CAPLUS

CN Benzeneacetamide, N-[(1R)-1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-3-(trifluoromethyl)- (9CI) (CA INDEX NAME)

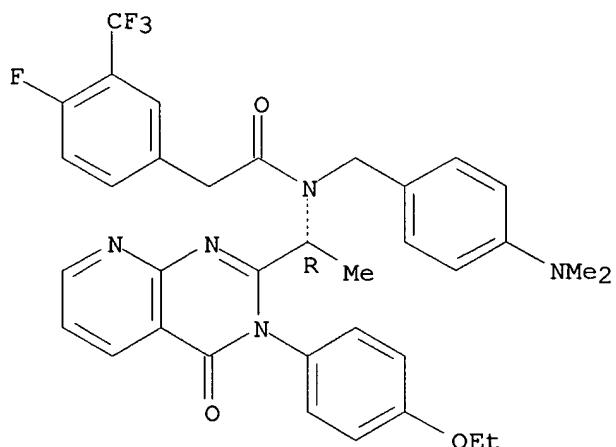
Absolute stereochemistry.



RN 473720-24-0 CAPLUS

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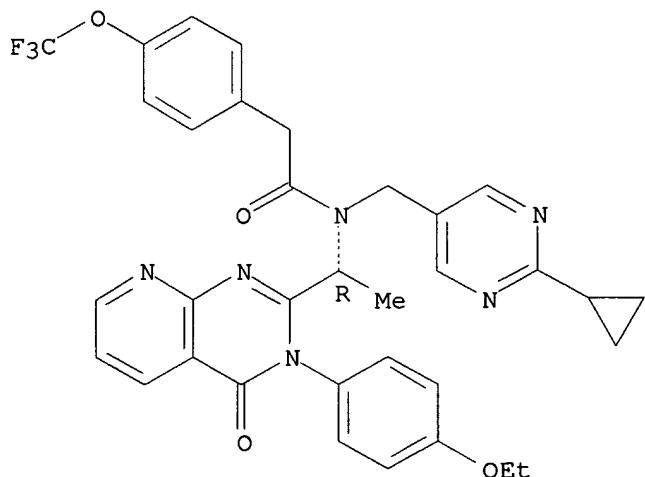
Absolute stereochemistry.



RN 473720-25-1 CAPLUS

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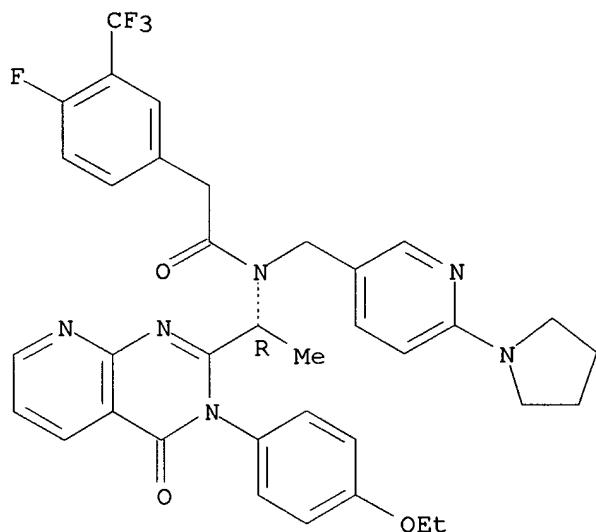
Absolute stereochemistry.



RN 473720-26-2 CAPLUS

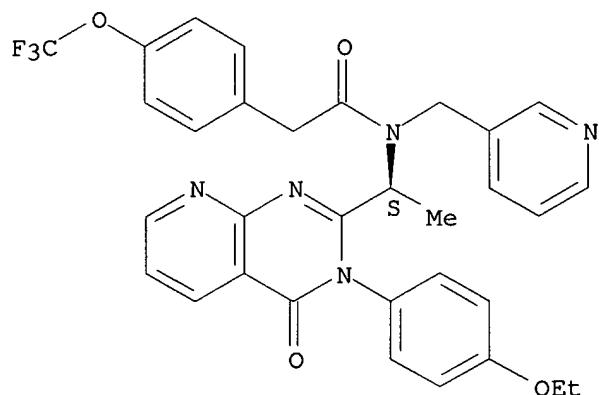
CN Benzeneacetamide, N-[(1R)-1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-N-[[6-(1-pyrrolidinyl)-3-pyridinyl]methyl]-3-(trifluoromethyl)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



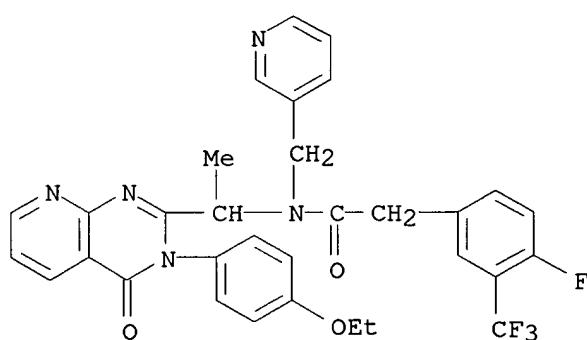
CN Benzeneacetamide, N-[(1S)-1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-N-(3-pyridinylmethyl)-4-(trifluoromethoxy)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



RN 473722-68-8 CAPLUS

CN Benzeneacetamide, N-[1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-4-fluoro-N-(3-pyridinylmethyl)-3-(trifluoromethyl)-(9CI) (CA INDEX NAME)



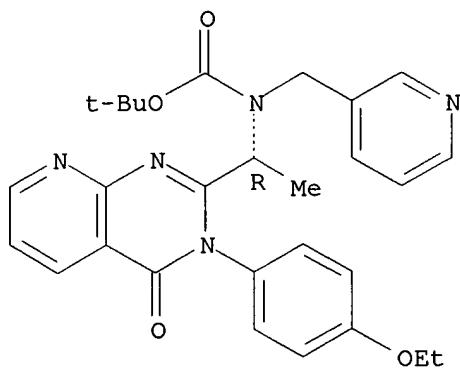
IT 473720-93-3P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(intermediate; preparation of N-(heteroarylalkyl)acylamides as CXCR3 antagonists for treatment of inflammatory or immune conditions)

RN 473720-93-3 CAPLUS

CN Carbamic acid, [(1R)-1-[3-(4-ethoxyphenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl](3-pyridinylmethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.



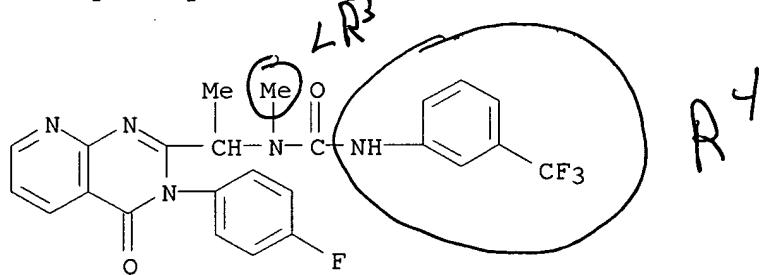
L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2003 ACS

2001:208250 Document No. 134:252352 Preparation of 3-aryl-2-arylureidoalkylquinazolin-4-ones and related compounds as mediators of hedgehog signaling pathways.. Baxter, Anthony David; Boyd, Edward Andrew; Guichert, Oivin M.; Price, Stephen; Rubin, Lee D. (Curis, Inc., USA). PCT Int. Appl. WO 2001019800 A2 20010322, 177 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 2000-US25461 20000915. PRIORITY: US 1999-PV154526 19990916; US 1999-PV159412 19991014; US 1999-PV162899 19991101.

| | PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|----|---|------|----------|-----------------|----------|
| PI | WO 2001019800 | A2 | 20010322 | WO 2000-US25461 | 20000915 |
| | WO 2001019800 | A3 | 20011206 | | |
| | WO 2001019800 | C2 | 20021003 | | |
| | W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG | | | | |
| | EP 1216234 | A2 | 20020626 | EP 2000-963551 | 20000915 |
| | R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL | | | | |
| | JP 2003509414 | T2 | 20030311 | JP 2001-523380 | 20000915 |
| | US 6545005 | B1 | 20030408 | US 2000-663835 | 20000915 |
| IT | 330796-36-6P | | | | |
| | RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of 3-aryl-2-arylureidoalkylquinazolin-4-ones and related compds. as mediators of hedgehog signaling pathways) | | | | |

RN 330796-36-6 CAPLUS

CN Urea, N-[1-[3-(4-fluorophenyl)-3,4-dihydro-4-oxopyrido[2,3-d]pyrimidin-2-yl]ethyl]-N-methyl-N'-[3-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



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LOGOFF? (Y)/N/HOLD:.

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

FULL ESTIMATED COST

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156.29

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